

IN THE CLAIMS

1. (Amended) A compound of formula E-C_a-R-C_b-A, wherein E is a therapeutic or diagnostic agent, R is a reactive group, C_b and C_a are optional first and second connecting respectively, and A is an affinity group comprising any molecule or part of a molecule possessing specific binding determinants for a target molecule having an affinity for human serum albumin, wherein affinity group A comprises a sequence of amino acid residues -O₁-O₂-X₁-X₂-B in which the amino acid residues are independently selected from the group of all twenty naturally occurring amino acids.

A³
2. (Amended) A compound according to claim 58, wherein amino acid residue O₁ is selected from the group consisting of phenylalanine, arginine, glutamine, tyrosine, glutamic acid and tryptophan; amino acid residue O₂ is selected from the group consisting of leucine, arginine, glutamic acid, tryptophan and phenylalanine; amino acid residue X₁ is selected from the group consisting of phenylalanine, tryptophan, methionine and tyrosine; amino acid residue X₂ is selected from the group consisting of serine, arginine and glutamic acid; and amino acid residue B is selected from the group consisting of serine, arginine and glutamic acid.

Please CANCEL Claim 3.

4. (Amended) A compound according to claim 58, wherein one of the five amino acid residues is an L amino acid residue and the other four amino acid residues are D amino acid residues.

A^d
5. (Amended) A compound according to claim 2, wherein the L-amino acid residue is selected from the group consisting of the amino acid residue O₂, the amino acid residue X₁, and the amino acid residue X₂.

6. (Amended) A compound according to claim 58, wherein one of the five amino acid residues is a D-amino acid residue and the other four amino acid residues are L-amino acid residues.

7. (Reiterated) A compound according to claim 6, wherein the D-amino acid residue is selected from the group consisting of the amino acid residue O₂, amino acid residue X₁, and amino acid residue X₂.

8. (Reiterated) A compound according to claim 7, wherein the D-amino acid residue is the amino acid residue O₂.

9. (Amended) A compound according to claim 58, wherein O₁ is phenylalanine and O₂ is leucine.

10. (Amended) A compound according to claim 58, wherein O₁ is arginine and O₂ is arginine.

11. (Amended) A compound according to claim 58, wherein O₁ is glutamine and O₂ is glutamic acid.

12. (Amended) A compound according to claim 58, wherein O₁ is glutamic acid and O₂ is tryptophan.

13. (Amended) A compound according to claim 58, wherein O₁ is tryptophan and O₂ is tryptophan.

14. (Amended) A compound according to claim 58, wherein O₁ is tryptophan and O₂ is glutamic acid.

15. (Amended) A compound according to claim 58, wherein X₁ is tyrosine.

16. (Amended) A compound according to claim 58, wherein X₂ is glutamic acid.

17. (Amended) A compound according to claim 58, wherein B is glutamic acid.

18. (Amended) A compound according to claim 58, wherein O₁ is phenylalanine, O₂ is D-leucine, X₁ is tyrosine, X₂ is glutamic acid, and B is glutamic acid.

19. (Amended) A compound according to claim 58, wherein the amino acid residue B is a C-terminal amino acid residue.

20. (Reiterated) A compound according to claim 19, wherein the affinity group comprises the amino acid sequence -O₁-O₂-X₁-X₂-B-NH₂.

21. (Amended) A compound according to claim 58, wherein the the reactive group comprises a functional group selected from the group consisting of carboxy, phosphoryl, alkyl esters, thioesters, phosphoesters, ortho esters, imidates, mixed anhydrides, amides, thioamine and disulphides.

22. (Amended) A compound according to claim 21, wherein C_b is absent and the reactive group is bonded directly to the O₁ amino acid residue in the affinity group.

23. (Reiterated) A compound according to claim 22, wherein the reactive group is bonded to the O₁ amino acid residue by an amide linkage.

24. (Amended) A compound according to claim 21, wherein the reactive group has the formula -X-R₁-C(O)-, wherein R₁ comprises a substituted or unsubstituted aromatic group and X is selected from the group consisting of S, O and N.

25. (Reiterated) A compound according to claim 24, wherein X is bonded directly to an aromatic carbon atom in R₁.

26. (Reiterated) A compound according to claim 24, wherein R₁ is unsubstituted phenyl.

27. (Amended) A compound according to claim 26, wherein -X- and -C(O)- are bonded to the phenyl in a para configuration.

46 cont

7

28. (Amended) A compound according to claim 24, wherein R₁ is phenyl substituted with one or more groups selected from the group consisting of a halogen, NO₂, SO₂NH₂, SO₂NHF, CF₃, CCl₃, CBr₃, C≡N, SO₃H, CO₂H, CHO, OH, NHCOCH₃, OCH₃, CH₃ and CH₂CH₃.

29. (Reiterated) A compound according to claim 24, wherein the reactive moiety is bonded directly to the O₁ residue via the carboxyl carbon.

30. (Amended) A compound according to claim 21 wherein C_b is present.

31. (Amended) A compound according to claim 28, wherein C_b is bonded to the reactive group via an ester, thioester, amide, sulfonate ester or sulfonamide linkage.

32. (Amended) A compound according to claim 30, wherein C_b is bonded to the O₁ amino acid residue in the affinity group via an ester, thioester, amide, sulfonamide, urea, thiourea or carbamate linkage.

33. (Amended) A compound according to claim 30, wherein C_b comprises a backbone chain of between about 1 and about 25 atoms.

34. (Amended) A compound according to claim 33, wherein C_b comprises a backbone chain of between about 2 and about 16 carbon atoms.

35. (Amended) A compound according to claim 30, wherein C_b comprises an unsaturated carbon atom backbone chain of between about 1 and about 25 atoms.

Please CANCEL Claims 36-39.

40. (Amended) A compound according to claim 58 wherein C_a is present.

41. (Amended) A compound according to claim 40, wherein C_a is bonded to E by an ester, thioester, amide, sulfonate ester or sulfonamide linkage.

42. (Amended) A compound according to claim 40, wherein C_a is bonded to the reactive group by an ester, thioester, amide or sulfonate ester linkage.

43. (Amended) A compound according to claim 40, wherein C_a comprises a backbone chain of between about 1 and about 25 atoms.

44. (Amended) A compound according to claim 43, wherein C_a comprises a backbone chain of between about 2 and about 16 carbon atoms.

45. (Amended) A compound according to claim 40, wherein C_a comprises an unsaturated carbon atom backbone chain of between about 1 and about 25 atoms.

46. (Amended) A compound according to claim 1, wherein the diagnostic agent comprises biotin.

47. (Amended) A compound according to claim 46, wherein biotin is bonded directly to the reactive group by an ester, thioester or amide linkage.

48. (Amended) A compound according to claim 46, wherein the reactive group has the formula -X-Ph-C(O)-, and wherein X is oxygen, sulfur or nitrogen.

49. (Amended) A compound according to claim 48, wherein the -X- and -C(O)- on the phenyl group are bonded in a para configuration.

50. (Amended) A compound according to claim 47 wherein C_a is present.

51. (Amended) A compound according to claim 50, wherein C_a is bonded to the biotin group by an amide linkage.